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NEWS 1		Web Page URLs for STN Seminar Schedule - N. America
NEWS 2	Apr 08	"Ask CAS" for self-help around the clock
NEWS 3	Apr 09	BEILSTEIN: Reload and Implementation of a New Subject Area
NEWS 4	Apr 09	ZDB will be removed from STN
NEWS 5	Apr 19	US Patent Applications available in IFICDB, IFIPAT, and IFIUDB
NEWS 6	Apr 22	Records from IP.com available in CAPLUS, HCAPLUS, and ZCAPLUS
NEWS 7	Apr 22	BIOSIS Gene Names now available in TOXCENTER
NEWS 8	Apr 22	Federal Research in Progress (FEDRIP) now available
NEWS 9	Jun 03	New e-mail delivery for search results now available
NEWS 10	Jun 10	MEDLINE Reload
NEWS 11	Jun 10	PCTFULL has been reloaded
NEWS 12	Jul 02	FOREGE no longer contains STANDARDS file segment
NEWS 13	Jul 22	USAN to be reloaded July 28, 2002; saved answer sets no longer valid
NEWS 14	Jul 29	Enhanced polymer searching in REGISTRY
NEWS 15	Jul 30	NETFIRST to be removed from STN
NEWS 16	Aug 08	CANCERLIT reload
NEWS 17	Aug 08	PHARMAMarketLetter(PHARMAML) - new on STN
NEWS 18	Aug 08	NTIS has been reloaded and enhanced
NEWS 19	Aug 19	Aquatic Toxicity Information Retrieval (AQUIRE) now available on STN
NEWS 20	Aug 19	IFIPAT, IFICDB, and IFIUDB have been reloaded
NEWS 21	Aug 19	The MEDLINE file segment of TOXCENTER has been reloaded
NEWS 22	Aug 26	Sequence searching in REGISTRY enhanced
NEWS 23	Sep 03	JAPIO has been reloaded and enhanced
NEWS 24	Sep 16	Experimental properties added to the REGISTRY file
NEWS 25	Sep 16	Indexing added to some pre-1967 records in CA/CAPLUS
NEWS 26	Sep 16	CA Section Thesaurus available in CAPLUS and CA
NEWS 27	Oct 01	CASREACT Enriched with Reactions from 1907 to 1985

NEWS EXPRESS	October 14 CURRENT WINDOWS VERSION IS V6.01, CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP), AND CURRENT DISCOVER FILE IS DATED 01 OCTOBER 2002
NEWS HOURS	STN Operating Hours Plus Help Desk Availability
NEWS INTER	General Internet Information
NEWS LOGIN	Welcome Banner and News Items
NEWS PHONE	Direct Dial and Telecommunication Network Access to STN
NEWS WWW	CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

* * * * * * * * * * * * * * * * STN Columbus * * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 17:14:20 ON 16 OCT 2002

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 17:14:29 ON 16 OCT 2002

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 15 OCT 2002 HIGHEST RN 461638-40-4
DICTIONARY FILE UPDATES: 15 OCT 2002 HIGHEST RN 461638-40-4

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=> e benzyl(octahydro-4,7-methanoinden-5-yl)amine/cn
E1 1 BENZYL(O-HYDROXYPHENYL)CARBINOL/CN
E2 1 BENZYL(O-METHOXY-.ALPHA.-PHENETHYLBENZYL)DIMETHYLAMMONIUM BR
OMIDE/CN
E3 0 --> BENZYL(OCTAHYDRO-4,7-METHANOINDEN-5-YL)AMINE/CN
E4 1 BENZYL(OCTYLOXYCARBONYLMETHYL)(OCTYLCARBAMOYLMETHYL)AMINE/CN
E5 1 BENZYL(P-(2-(BENZYLMETHYLAMINO)ETHOXY)-.BETA.-HYDROXY-.ALPHA
.-METHYLPHENETHYL)DIMETHYLAMMONIUM BROMIDE METHOBROMIDE/CN
E6 1 BENZYL(P-(2-CHLORO-4-DIETHYLAMINOPHENYLAZO)PHENACYL)DIMETHYL
AMMONIUM CHLORIDE/CN
E7 1 BENZYL(P-(3,5-DICARBOXY-1,4-DIHYDRO-2,6-DIMETHYL-4-PYRIDYL)P

```
                    HENYL)DIMETHYLAMMONIUM IODIDE, DIMETHYL ESTER/CN
E8            1
BENZYL(P-(3-HYDROXY-1-OXOINDEN-2-YL)PHENYL)DIMETHYLAMMONIUM
                    HYDROXIDE, INNER SALT/CN
E9            1     BENZYL(P-(DIMETHYLAMINO)PHENYL)DIMETHYLPHOSPHONIUM
CHLORIDE/
                    CN
E10           1     BENZYL(P-(DIMETHYLAMINO)PHENYL)DIPHENYLPHOSPHONIUM
BROMIDE/C
                    N
E11           1     BENZYL(P-(DIMETHYLAMINO)PHENYL)DIPHENYLPHOSPHONIUM
CHLORIDE/
                    CN
E12           1     BENZYL(P-1-BUTENYLPHENYL)DIMETHYLAMMONIUM BROMIDE/CN

=> e benzyl(octahydro-4,7-methanoinden-5-yl)amine
E1            1     BENZYILDENE/BI
E2       205225     BENZYL/BI
E3            0 --> BENZYL(OCTAHYDRO-4,7-METHANOINDEN-5-YL)AMINE/BI
E4            3     BENZYLA/BI
E5            1     BENZYLABRES/BI
E6            1     BENZYLABRESOL/BI
E7            1     BENZYLABRESOLINE/BI
E8            2     BENZYLACENAPHTH/BI
E9            2     BENZYLACENAPHTHO/BI
E10         193     BENZYLACET/BI
E11           3     BENZYLACETALDEHYDE/BI
E12          22     BENZYLACETAMIDE/BI

=> file caplus
COST IN U.S. DOLLARS                          SINCE FILE      TOTAL
                                                   ENTRY    SESSION
FULL ESTIMATED COST                                1.52       1.73

FILE 'CAPLUS' ENTERED AT 17:16:57 ON 16 OCT 2002
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FILE COVERS 1907 - 16 Oct 2002  VOL 137 ISS 16
FILE LAST UPDATED: 15 Oct 2002  (20021015/ED)
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This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

```
=> e benzyl(octahydro-4,7-methanoinden-5-yl)amine
E1            1      BENZYKPIPERIDONE/BI
E2       134346      BENZYL/BI
E3            0 --> BENZYL(OCTAHYDRO-4,7-METHANOINDEN-5-YL)AMINE/BI
E4            1      BENZYL0/BI
E5            2      BENZYL1/BI
E6            4      BENZYL2/BI
E7            6      BENZYL3/BI
E8            1      BENZYL3SN/BI
E9            4      BENZYL4/BI
E10           1      BENZYL8/BI
E11           2      BENZYLA/BI
E12           1      BENZYLABIETAMIDE/BI

=> e benzyl(octahydro-4,7-methanoinden-5-yl)amine/cn
   REG1stRY INITIATED
Substance data EXPAND from CAS REGISTRY in progress...

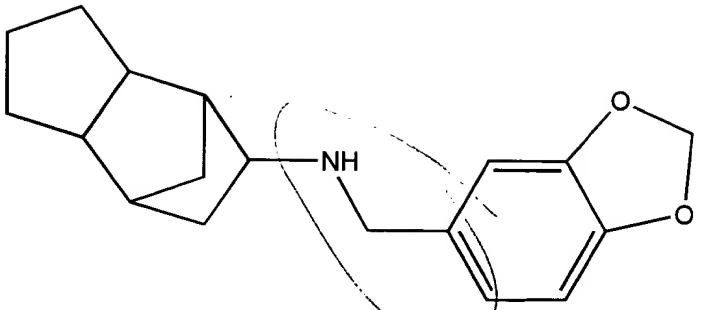


E1            1      BENZYL(O-HYDROXYPHENYL)CARBINOL/CN
E2            1      BENZYL(O-METHOXY-.ALPHA.-PHENETHYLBENZYL)DIMETHYLAMMONIUM
BR
                     OMIDE/CN
E3            0 --> BENZYL(OCTAHYDRO-4,7-METHANOINDEN-5-YL)AMINE/CN
E4            1
BENZYL(OCTYLOXYCARBONYLMETHYL)(OCTYLCARBAMOYLMETHYL)AMINE/CN
E5            1
BENZYL(P-(2-(BENZYLMETHYLAMINO)ETHOXY)-.BETA.-HYDROXY-.ALPHA
                     .-METHYLPHENETHYL)DIMETHYLAMMONIUM BROMIDE METHOBROMIDE/CN
E6            1
BENZYL(P-(2-CHLORO-4-DIETHYLAMINOPHENYLAZO)PHENACYL)DIMETHYL
                     AMMONIUM CHLORIDE/CN
E7            1
BENZYL(P-(3,5-DICARBOXY-1,4-DIHYDRO-2,6-DIMETHYL-4-PYRIDYL)P
                     HENYL)DIMETHYLAMMONIUM IODIDE, DIMETHYL ESTER/CN
E8            1
BENZYL(P-(3-HYDROXY-1-OXOINDEN-2-YL)PHENYL)DIMETHYLAMMONIUM
                     HYDROXIDE, INNER SALT/CN
E9            1      BENZYL(P-(DIMETHYLAMINO)PHENYL)DIMETHYLPHOSPHONIUM
CHLORIDE/
                     CN
E10           1      BENZYL(P-(DIMETHYLAMINO)PHENYL)DIPHENYLPHOSPHONIUM
BROMIDE/C
                     N
E11           1      BENZYL(P-(DIMETHYLAMINO)PHENYL)DIPHENYLPHOSPHONIUM
CHLORIDE/
                     CN
E12           1      BENZYL(P-1-BUTENYLPHENYL)DIMETHYLAMMONIUM BROMIDE/CN
```

amide

564/163

164



amide

384

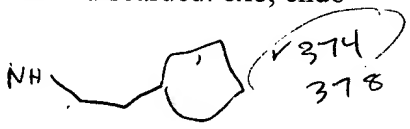
387 polycyclic

391 unsub

389 OH to ring

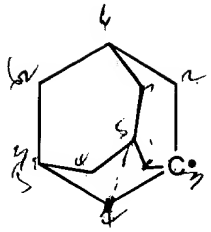
exo/endo-benzo[1,3]dioxol-5-ylmethyl(octahydro-4,7-methanoinden-5-yl)amine

Caution: Stereochemical terms discarded: exo, endo

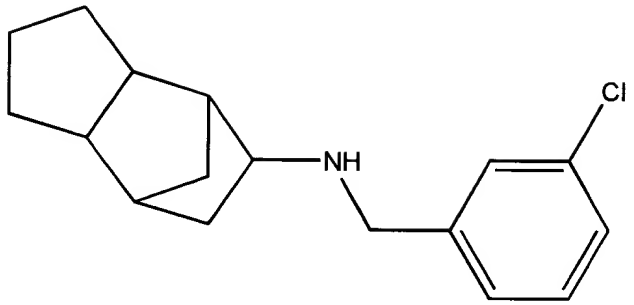


374

378 naphthyl



adamantyl



exo/endo-(3chlorobenzyl)(octahydro-4,7-methanoinden-5-yl)amine

Caution: Stereochemical terms discarded: exo, endo

562/442

558/422

560/37

564/163